

Discussion of “Feature Matching in Time Series Modeling” by Y. Xia and H. Tong

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Many congratulations to Professors Xia and Tong for another stimulating paper initiated from their own creative thinking. The base point of the proposed approach is the fact that most, if not all, statistical models are wrong. This not only applies to time series models, as a statistical model is, hopefully, a simplified representation of the truth. At the best it catches some features of the unknown underlying population. While the understanding of this nature is within the common wisdom, most statistical inference methods are confined to the framework which assumes that the true model is a member of the family of models concerned. The approach advocated in this paper acknowledges explicitly that the assumed model is not the truth, and indeed it is advantageous sometimes not to read too much into the assumed model. For example, the authors have articulated elegantly that if our interest lies in catching the linear dynamical structure, we should not use the (Gaussian) maximum likelihood estimation which effectively minimizes the one-step-ahead prediction errors only, and in fact a better fitted autocovariance is resulted from minimizing up to m -step-ahead predictions for $m > 1$.

Following the lead of the authors, it seems to make sense to take on board the concern for “wrong models” at the stage of the model selection, too, as hinted at the end of the paper. In the way, this has been actively researched in the context of model selection. However, a difference here is to use a different measure for “goodness of fit” instead of likelihood (or log-likelihood). Let us consider a simple case: fit a linear AR(p) model to observations y_1, \dots, y_n

from a stationary time series with mean 0, where the order p is to be determined by the data, too. Let $\mathbf{y}_{t,p} = (y_t, y_{t-1}, \dots, y_{t-p+1})'$. Based on an AR(p) model (with independent innovations), the best predictor at the time t for a future value y_{t+m} should be a linear combination of the p components of $\mathbf{y}_{t,p}$. In fact the best linear predictor based on $\mathbf{y}_{t,p}$ is $\boldsymbol{\alpha}'_{m,p} \mathbf{y}_{t,p}$ with

$$(1) \quad \begin{aligned} \boldsymbol{\alpha}_{m,p} &= \boldsymbol{\Gamma}_p^{-1} \boldsymbol{\gamma}_{m,p} \\ &= \arg \min_{\boldsymbol{\gamma}} E\{(y_{t+m} - \boldsymbol{\alpha}' \mathbf{y}_{t,p})^2\}, \end{aligned}$$

where $\boldsymbol{\Gamma}_p$ is a $p \times p$ matrix with $\gamma(j-i)$ as its (i, j) th element, $\boldsymbol{\gamma}_{m,p}$ is a $p \times 1$ vector with $\gamma(m+i-1)$ as its i th element, and $\gamma(\cdot)$ denotes the autocovariance function of y_t . In fact (1) holds for any stationary process. However, if we fit y_t with an AR(p), its autocovariance function $\gamma(\cdot)$ is then determined by $\boldsymbol{\theta}_p$ —the parameters in an AR(p) model. Put $\boldsymbol{\alpha}_{m,p} = \boldsymbol{\alpha}_{m,p}(\boldsymbol{\theta}_p)$. Then $\mathbf{y}'_{t,p} \boldsymbol{\alpha}_{m,p}(\boldsymbol{\theta}_p)$ is the best predictor for y_{t+m} based on an AR(p) model. Using the “matching up-to- m -step-ahead point predictions” approach of Section 2.1, we estimate $\boldsymbol{\theta}_p$ (for p given) by

$$\hat{\boldsymbol{\theta}}_p = \arg \min_{\boldsymbol{\theta}_p} Q_p(\boldsymbol{\theta}_p),$$

where

$$Q_p(\boldsymbol{\theta}_p) = \frac{1}{m} \sum_{k=1}^m \frac{1}{n-k-p+1} \cdot \sum_{t=p}^{n-k} \{y_{t+k} - \mathbf{y}'_{t,p} \boldsymbol{\alpha}_{k,p}(\boldsymbol{\theta}_p)\}^2.$$

However, we cannot choose p by minimizing $Q_p(\hat{\boldsymbol{\theta}}_p)$, as $Q_p(\hat{\boldsymbol{\theta}}_p)$ is likely to decrease as p increases.

To appreciate the difficulties involved, let us first consider the “ideal world” where the (true) distribution of $\{y_t\}$ is known. Then we should estimate $\boldsymbol{\theta}_p$ by

$$\tilde{\boldsymbol{\theta}}_p = \arg \min_{\boldsymbol{\theta}_p} Q_p^*(\boldsymbol{\theta}_p),$$

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This is an electronic reprint of the original article published by the [Institute of Mathematical Statistics](#) in *Statistical Science*, 2011, Vol. 26, No. 1, 57–58. This reprint differs from the original in pagination and typographic detail.

where

$$Q_p^*(\boldsymbol{\theta}_p) = \frac{1}{m} \sum_{k=1}^m E[\{y_{t+k} - \mathbf{y}'_{t,p} \boldsymbol{\alpha}_{k,p}(\boldsymbol{\theta}_p)\}^2].$$

Unfortunately $Q_p^*(\tilde{\boldsymbol{\theta}}_p)$ still decreases as p increases. The information (e.g., the variance) of the noise component of y_t is required in order to know when to stop. This is the standard problem in model selection even for linear regression. One way to get away from this requirement is to take the log-transformation. Namely, we define

$$\begin{aligned} L^*(p) &= \log\{Q_p^*(\tilde{\boldsymbol{\theta}}_p)\} \\ &= \log\left\{\frac{1}{m} \sum_{k=1}^m E[\{y_{t+k} - \mathbf{y}'_{t,p} \boldsymbol{\alpha}_{k,p}(\tilde{\boldsymbol{\theta}}_p)\}^2]\right\}. \end{aligned}$$

When p is in the range on which $Q_p^*(\tilde{\boldsymbol{\theta}}_p)$ varies slowly (with respect to p), it holds that

$$L^*(p) - L^*(p+1) \approx \frac{Q_p^*(\tilde{\boldsymbol{\theta}}_p) - Q_{p+1}^*(\tilde{\boldsymbol{\theta}}_{p+1})}{Q_{p+1}^*(\tilde{\boldsymbol{\theta}}_{p+1})}.$$

Intuitively we would like to choose the smallest p such that the decrease $L^*(p) - L^*(p+1)$ is smaller than an appropriate but unknown constant. In practice, we may use $L(p) \equiv \log\{Q_p(\hat{\boldsymbol{\theta}}_p)\}$ to replace $L^*(p)$,

and choose p to minimize

$$L(p) + E\{L^*(p) - L(p)\}.$$

This is in the same spirit of AIC in the sense that the bias $E\{L^*(p) - L(p)\}$ serves as the penalty for the model complexity. When the true model of y_t is not AR, this bias does not admit a simple asymptotic expression such as AIC even when $m = 1$; see, for example, Konishi and Kitagawa (1996). One may also consider to develop some resampling estimates for this bias.

The above line of thinking is provoked from reading this interesting paper which will serve as an inspiration for further research in tackling the issues related to the lack of a true model. Then one may quibble over the use of the phrase “catch-all approach.” If a model could catch all the features, it should be the true model, or at least pragmatically so. One message from the paper is that one should fit (and perhaps also choose) a model according to a specified purpose in hand, and a good statistical modeling is to catch the features of interest for a particular purpose.

REFERENCES

- KONISHI, S. and KITAGAWA, G. (1996). Generalised information criteria in model selection. *Biometrika* **83** 875–890. [MR1440051](#)